

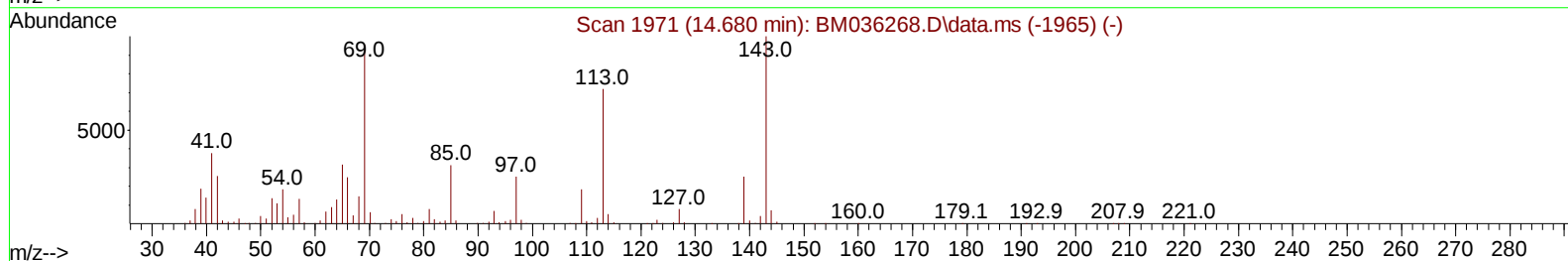
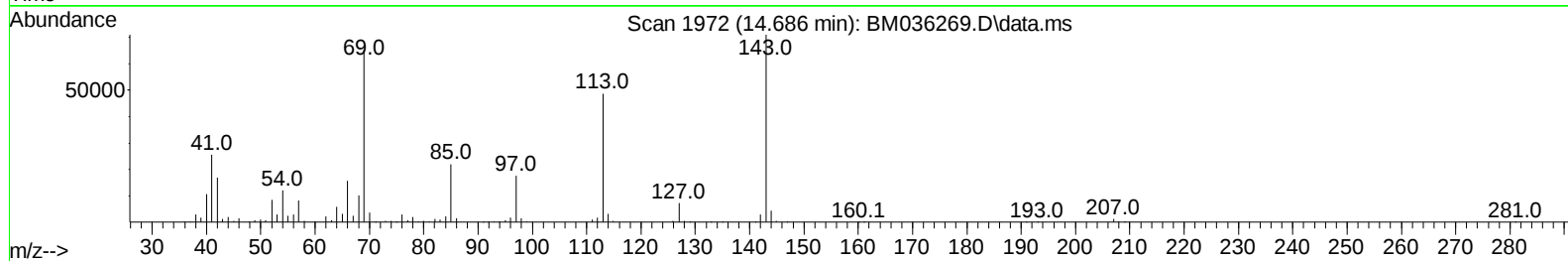
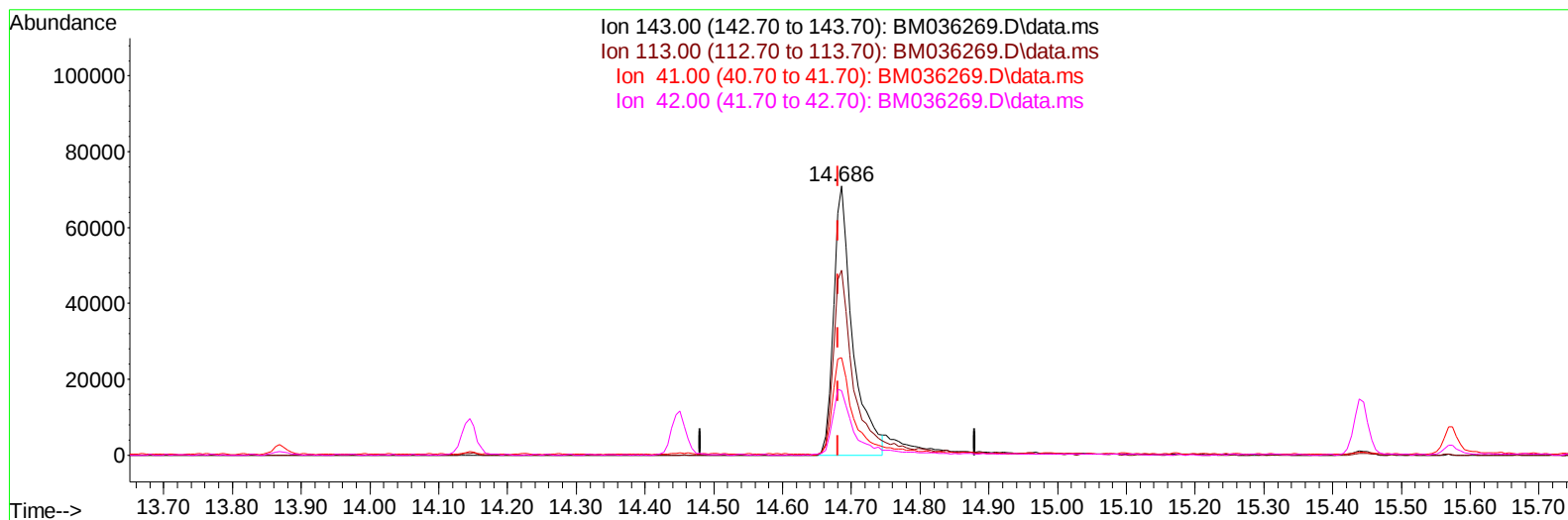
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081422\
Data File : BM036269.D
Acq On : 14 Aug 2022 14:18
Operator : CG/JU
Sample : PB146982BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK982

Manual IntegrationsAPPROVED

Quant Time: Aug 14 23:36:56 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Aug 14 23:34:05 2022
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/15/2022
Supervised By :mohammad ahmed 08/16/2022



TIC: BM036269.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.686min (+ 0.006) 15.75 ng/u1

response 136426

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	68.70
41.00	41.00	36.22
42.00	29.30	23.96

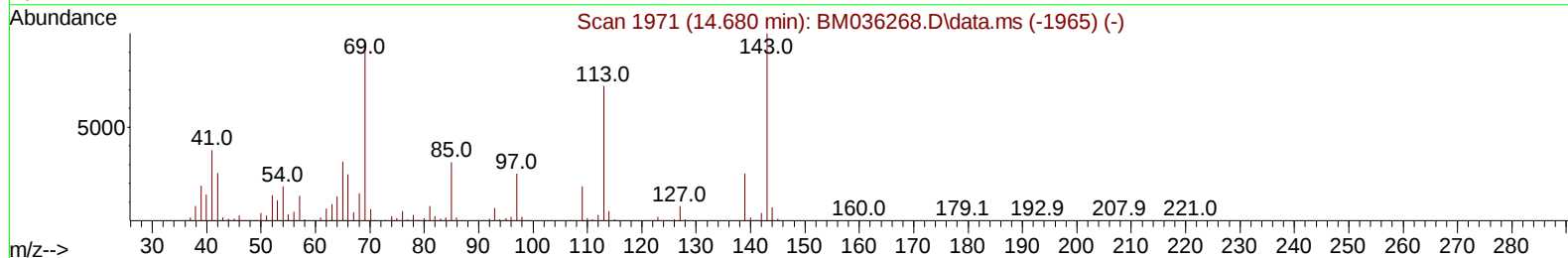
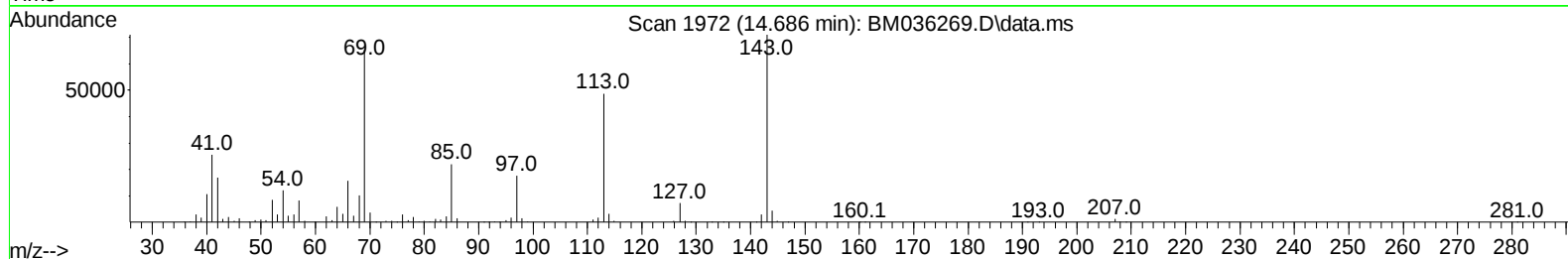
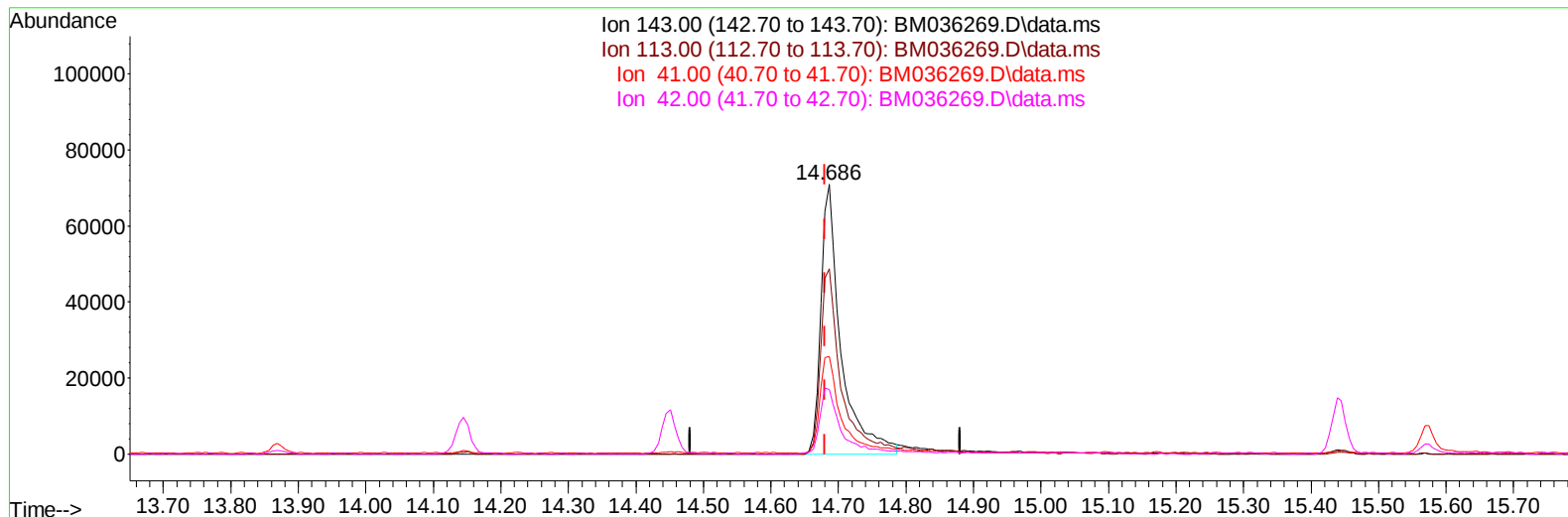
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TIC: BM036269.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.686min (+ 0.006) 16.78 ng/ul m

response 145347

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	68.70
41.00	41.00	36.22
42.00	29.30	23.96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	238349	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	1046934	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164	619059	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	1212341	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	1086319	20.000	ng/ul	0.00
88) Perylene-d12	23.709	264	1107122	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	27895	4.404	ng/uL	0.00
4) Pyridine-d5	3.640	84	186658	10.757	ng/ul	0.00
7) Phenol-d5	6.992	99	456423	22.835	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.145	67	303253	22.759	ng/ul	0.00
11) 2-Chlorophenol-d4	7.339	132	387308	23.345	ng/ul	0.00
15) 4-Methylphenol-d8	8.533	113	358863	22.162	ng/ul	0.00
21) Nitrobenzene-d5	8.980	128	188999	22.782	ng/ul	0.00
24) 2-Nitrophenol-d4	9.698	143	197605	24.448	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.239	165	329829	22.563	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	294725	11.909	ng/ul	0.00
46) Dimethylphthalate-d6	13.868	166	1051465	24.935	ng/ul	0.00
49) Acenaphthylene-d8	14.145	160	1225566	22.779	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	145347m	16.785	ng/ul	0.00
60) Fluorene-d10	15.445	176	925816	24.296	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	124128	19.048	ng/ul	0.00
73) Anthracene-d10	17.292	188	1443173	26.330	ng/ul	0.00
81) Pyrene-d10	19.586	212	1565990	26.011	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.556	264	1458899	25.956	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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